



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 29, 2024 – 04:18 PM JST

PDB ID : 8JT1
Title : COLLAGENASE FROM GRIMONTIA (VIBRIO) HOLLISAE 1706B COM-
PLEXED WITH GLY-PRO-HYP-GLY-PRO-HYP
Authors : Ueshima, S.; Yaskawa, K.; Takita, T.; Mikami, B.
Deposited on : 2023-06-21
Resolution : 2.00 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.36
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36

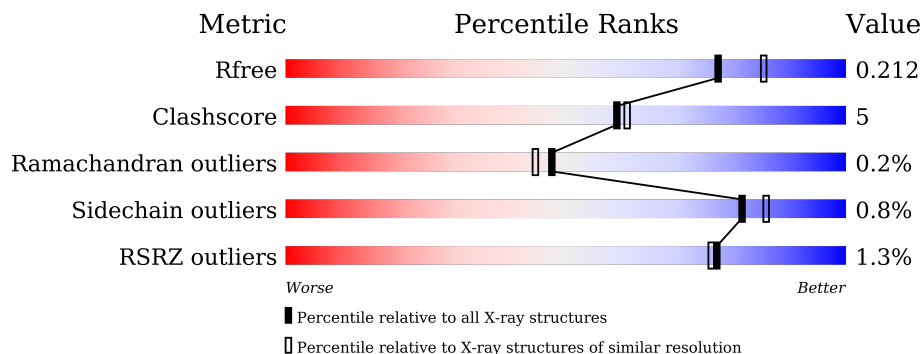
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	559	
1	B	559	
2	C	6	
2	E	6	
2	F	6	
2	G	6	

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Mol	Chain	Length	Quality of chain
2	I	6	
2	J	6	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	EDO	B	732	-	-	X	-
5	EDO	B	742	-	-	X	-

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 9660 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Microbial collagenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	532	4251	2692	713	827	19	0	9	0
1	B	532	4246	2689	711	827	19	0	7	0

- Molecule 2 is a protein called 6-mer peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	6	39	24	6	9	0	0	0
2	E	6	39	24	6	9	0	0	0
2	F	3	19	12	3	4	0	0	0
2	G	6	39	24	6	9	0	0	0
2	I	6	39	24	6	9	0	0	0
2	J	3	19	12	3	4	0	0	0

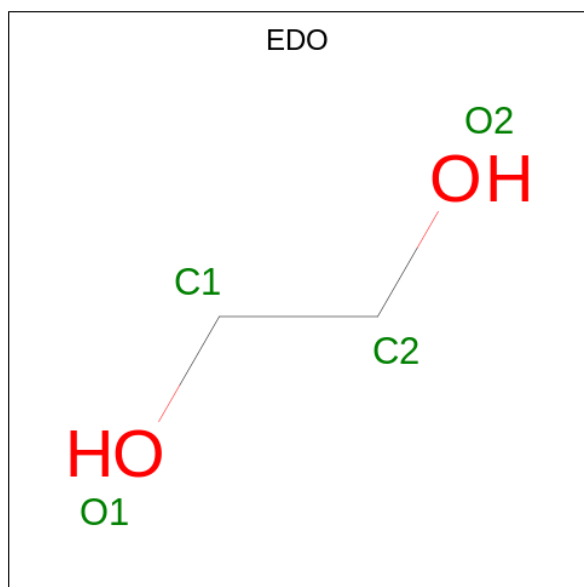
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Zn	0	0
			1	1		
3	B	1	Total	Zn	0	0
			1	1		

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	4	Total Ca 4 4	0	0
4	B	4	Total Ca 4 4	0	0

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0
5	A	1	Total C O 4 2 2	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	A	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0

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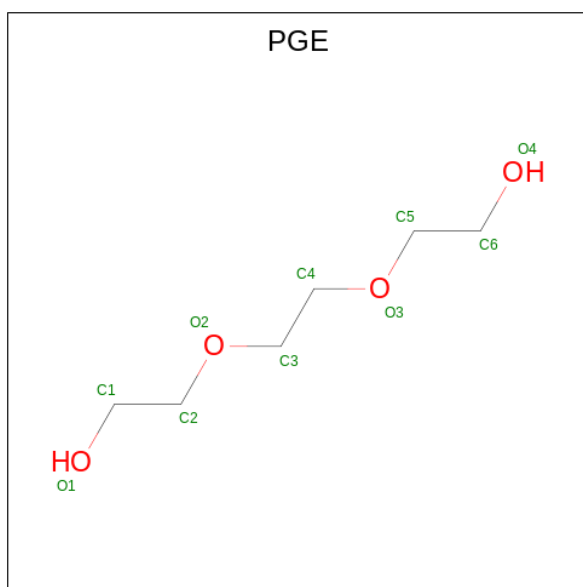
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0
5	B	1	Total 4	C 2	O 2	0	0

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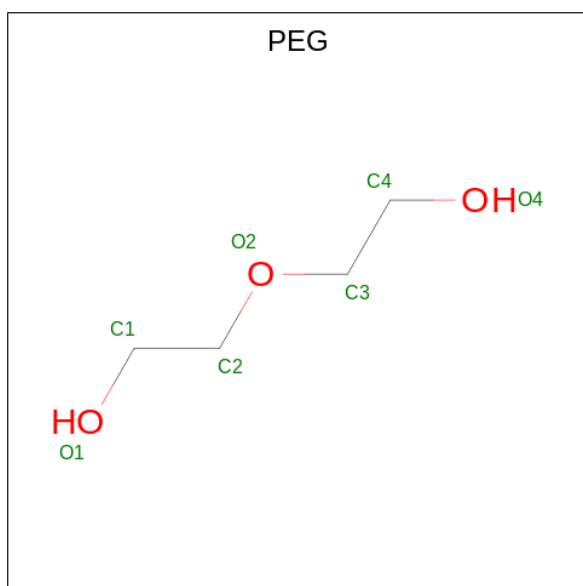
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	1
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is TRIETHYLENE GLYCOL (three-letter code: PGE) (formula: C₆H₁₄O₄) (labeled as "Ligand of Interest" by depositor).



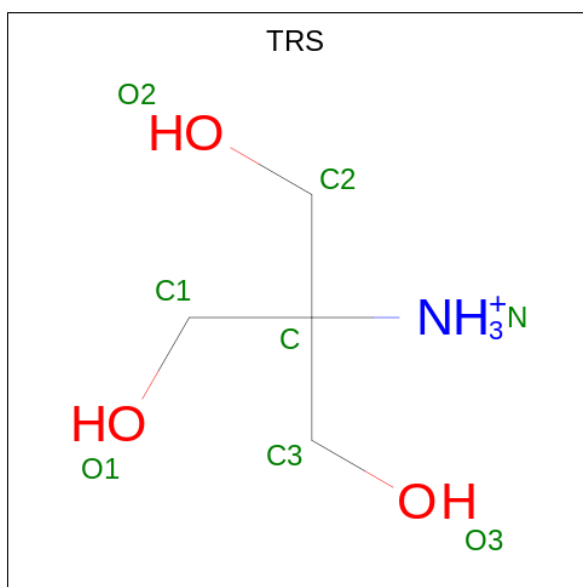
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 8 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	N			O
8	B	1	8	4	1	3	0	0

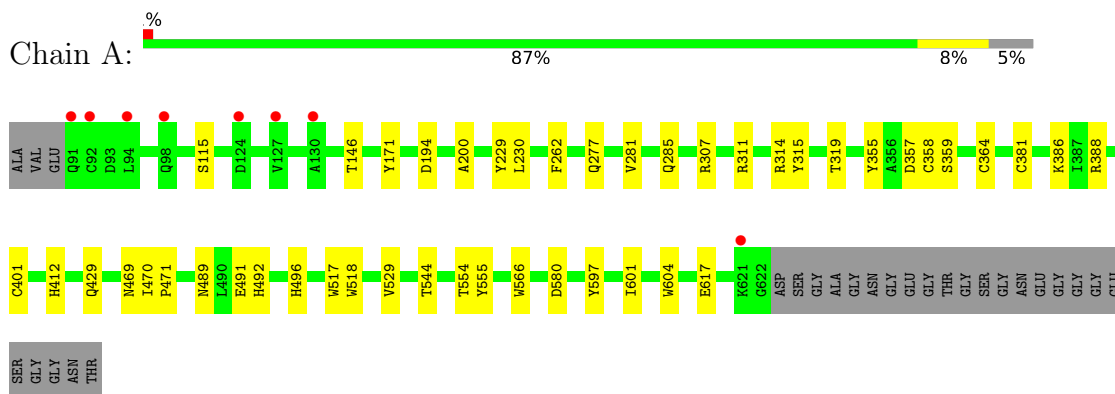
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	269	Total	O	0	2
			271	271		
9	B	288	Total	O	0	0
			288	288		
9	C	5	Total	O	0	0
			5	5		
9	E	7	Total	O	0	0
			7	7		
9	G	2	Total	O	0	0
			2	2		
9	I	4	Total	O	0	0
			4	4		
9	J	1	Total	O	0	0
			1	1		

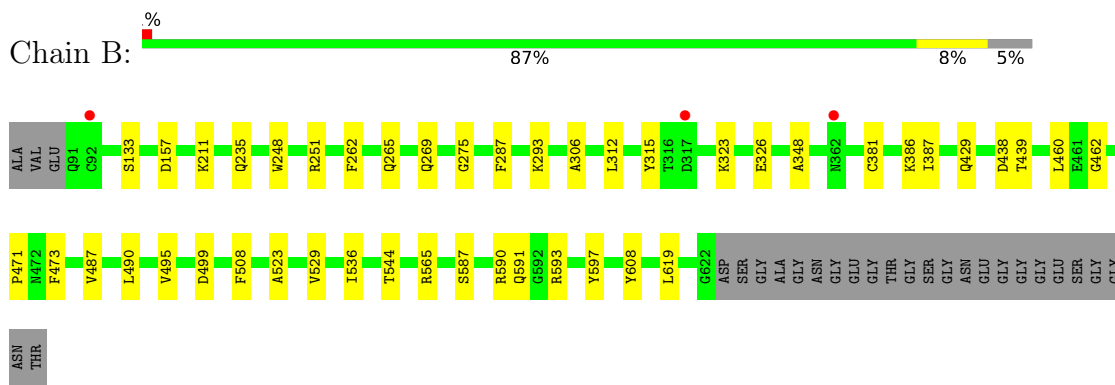
3 Residue-property plots i

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

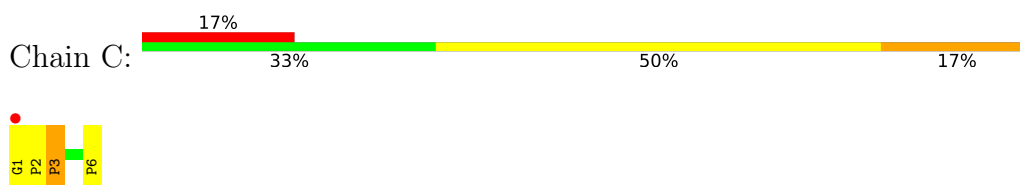
- Molecule 1: Microbial collagenase



- Molecule 1: Microbial collagenase



- Molecule 2: 6-mer peptide

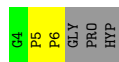


- Molecule 2: 6-mer peptide

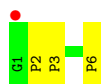




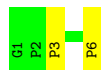
- Molecule 2: 6-mer peptide



- Molecule 2: 6-mer peptide



- Molecule 2: 6-mer peptide



- Molecule 2: 6-mer peptide



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	178.02Å 74.89Å 135.13Å 90.00° 126.02° 90.00°	Depositor
Resolution (Å)	37.44 – 2.00 46.03 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.6 (37.44-2.00) 99.6 (46.03-2.00)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.34 (at 2.00Å)	Xtrriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.176 , 0.214 0.176 , 0.212	Depositor DCC
R_{free} test set	4826 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	32.7	Xtrriage
Anisotropy	0.368	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 51.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9660	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.63% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PEG, TRS, HYP, EDO, CA, ZN, PGE

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.37	0/4386	0.55	0/5977
1	B	0.37	0/4373	0.55	0/5960
2	C	0.28	0/22	0.54	0/28
2	E	0.30	0/22	0.55	0/28
2	F	0.22	0/11	0.63	0/14
2	G	0.32	0/22	0.44	0/28
2	I	0.26	0/22	0.44	0/28
2	J	0.31	0/11	0.63	0/14
All	All	0.37	0/8869	0.55	0/12077

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4251	0	3947	34	0
1	B	4246	0	3925	42	0
2	C	39	0	36	4	0
2	E	39	0	36	1	0
2	F	19	0	16	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	39	0	36	2	0
2	I	39	0	36	0	0
2	J	19	0	16	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	4	0	0	0	0
4	B	4	0	0	0	0
5	A	176	0	264	21	0
5	B	180	0	270	26	0
6	A	10	0	14	0	0
7	B	7	0	10	2	0
8	B	8	0	12	0	0
9	A	271	0	0	3	0
9	B	288	0	0	3	0
9	C	5	0	0	2	0
9	E	7	0	0	0	0
9	G	2	0	0	0	0
9	I	4	0	0	0	0
9	J	1	0	0	0	0
All	All	9660	0	8618	84	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 5.

The worst 5 of 84 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:554:THR:H	5:A:727:EDO:H21	1.35	0.91
1:B:251:ARG:HE	7:B:751:PEG:H22	1.33	0.91
1:B:544:THR:H	5:B:709:EDO:H11	1.45	0.82
1:B:462:GLY:H	5:B:732:EDO:H12	1.51	0.76
1:B:133:SER:HB2	5:B:723:EDO:H21	1.70	0.72

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	539/559 (96%)	534 (99%)	4 (1%)	1 (0%)	47	44
1	B	537/559 (96%)	528 (98%)	8 (2%)	1 (0%)	47	44
2	C	3/6 (50%)	3 (100%)	0	0	100	100
2	E	3/6 (50%)	3 (100%)	0	0	100	100
2	F	1/6 (17%)	1 (100%)	0	0	100	100
2	G	3/6 (50%)	3 (100%)	0	0	100	100
2	I	3/6 (50%)	3 (100%)	0	0	100	100
2	J	1/6 (17%)	1 (100%)	0	0	100	100
All	All	1090/1154 (94%)	1076 (99%)	12 (1%)	2 (0%)	47	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	529	VAL
1	B	529	VAL

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	437/442 (99%)	433 (99%)	4 (1%)	78	83
1	B	434/442 (98%)	430 (99%)	4 (1%)	78	83
2	C	2/2 (100%)	2 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	2/2 (100%)	2 (100%)	0	100	100
2	F	1/2 (50%)	1 (100%)	0	100	100
2	G	2/2 (100%)	2 (100%)	0	100	100
2	I	2/2 (100%)	2 (100%)	0	100	100
2	J	1/2 (50%)	1 (100%)	0	100	100
All	All	881/896 (98%)	873 (99%)	8 (1%)	81	83

5 of 8 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	597	TYR
1	B	473	PHE
1	B	326[A]	GLU
1	A	617	GLU
1	B	326[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	96	GLN
1	A	135	ASN
1	A	606	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

10 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HYP	C	6	2	9,9,9	1.26	1 (11%)	10,12,12	1.60	2 (20%)
2	HYP	C	3	2,3	6,8,9	0.93	0	5,10,12	1.54	1 (20%)
2	HYP	E	3	2	6,8,9	0.98	0	5,10,12	1.57	2 (40%)
2	HYP	I	6	2	9,9,9	1.01	0	10,12,12	1.15	1 (10%)
2	HYP	J	6	2	6,8,9	1.11	0	5,10,12	2.73	3 (60%)
2	HYP	E	6	2	9,9,9	1.15	1 (11%)	10,12,12	1.33	3 (30%)
2	HYP	G	3	2,3	6,8,9	1.07	0	5,10,12	1.57	1 (20%)
2	HYP	I	3	2	6,8,9	0.92	0	5,10,12	1.41	1 (20%)
2	HYP	F	6	2	6,8,9	0.92	0	5,10,12	2.20	2 (40%)
2	HYP	G	6	2	9,9,9	1.05	0	10,12,12	1.63	1 (10%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HYP	C	6	2	-	1/4/13/13	0/1/1/1
2	HYP	C	3	2,3	-	0/0/11/13	0/1/1/1
2	HYP	E	3	2	-	0/0/11/13	0/1/1/1
2	HYP	I	6	2	-	4/4/13/13	0/1/1/1
2	HYP	J	6	2	-	0/0/11/13	0/1/1/1
2	HYP	E	6	2	-	1/4/13/13	0/1/1/1
2	HYP	G	3	2,3	-	0/0/11/13	0/1/1/1
2	HYP	I	3	2	-	0/0/11/13	0/1/1/1
2	HYP	F	6	2	-	0/0/11/13	0/1/1/1
2	HYP	G	6	2	-	1/4/13/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	6	HYP	CA-C	2.50	1.59	1.52
2	C	6	HYP	CA-C	2.25	1.58	1.52

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	6	HYP	O-C-CA	-4.07	114.12	124.78
2	F	6	HYP	CB-CG-CD	-3.37	99.13	103.27

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	6	HYP	CB-CG-CD	-3.02	99.56	103.27
2	G	6	HYP	O-C-CA	-2.72	113.32	122.26
2	E	6	HYP	CB-CG-CD	-2.51	100.18	103.27

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	6	HYP	OXT-C-CA-N
2	I	6	HYP	O-C-CA-N
2	I	6	HYP	O-C-CA-CB
2	C	6	HYP	OXT-C-CA-N
2	G	6	HYP	OXT-C-CA-N

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	3	HYP	1	0

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 102 ligands modelled in this entry, 10 are monoatomic - leaving 92 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	EDO	A	732	-	3,3,3	0.42	0	2,2,2	0.43	0
5	EDO	A	723	-	3,3,3	0.42	0	2,2,2	0.44	0
5	EDO	A	747	-	3,3,3	0.51	0	2,2,2	0.33	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	A	724	-	3,3,3	0.40	0	2,2,2	0.51	0
5	EDO	B	740	-	3,3,3	0.58	0	2,2,2	0.34	0
5	EDO	B	712	-	3,3,3	0.53	0	2,2,2	0.29	0
5	EDO	B	738	-	3,3,3	0.43	0	2,2,2	0.38	0
5	EDO	A	707	-	3,3,3	0.48	0	2,2,2	0.26	0
5	EDO	A	708	-	3,3,3	0.46	0	2,2,2	0.39	0
5	EDO	B	731	-	3,3,3	0.49	0	2,2,2	0.42	0
5	EDO	A	735	-	3,3,3	0.57	0	2,2,2	0.21	0
5	EDO	B	732	-	3,3,3	0.39	0	2,2,2	0.51	0
5	EDO	B	748	-	3,3,3	0.36	0	2,2,2	0.86	0
5	EDO	B	724	-	3,3,3	0.41	0	2,2,2	0.43	0
5	EDO	A	748	-	3,3,3	0.55	0	2,2,2	0.33	0
5	EDO	B	717	-	3,3,3	0.51	0	2,2,2	0.15	0
5	EDO	B	726	-	3,3,3	0.47	0	2,2,2	0.52	0
5	EDO	B	707	-	3,3,3	0.49	0	2,2,2	0.45	0
5	EDO	A	718	-	3,3,3	0.48	0	2,2,2	0.26	0
5	EDO	B	746	-	3,3,3	0.42	0	2,2,2	0.46	0
5	EDO	A	739	-	3,3,3	0.50	0	2,2,2	0.07	0
5	EDO	B	725	-	3,3,3	0.53	0	2,2,2	0.28	0
5	EDO	A	714	-	3,3,3	0.54	0	2,2,2	0.28	0
5	EDO	A	733	-	3,3,3	0.53	0	2,2,2	0.18	0
5	EDO	A	749	-	3,3,3	0.57	0	2,2,2	0.09	0
5	EDO	A	706	-	3,3,3	0.47	0	2,2,2	0.37	0
5	EDO	A	731	-	3,3,3	0.58	0	2,2,2	0.09	0
5	EDO	B	723	-	3,3,3	0.50	0	2,2,2	0.31	0
5	EDO	B	710	-	3,3,3	0.53	0	2,2,2	0.19	0
5	EDO	B	714	-	3,3,3	0.64	0	2,2,2	0.06	0
5	EDO	B	727	-	3,3,3	0.49	0	2,2,2	0.21	0
5	EDO	A	730	-	3,3,3	0.49	0	2,2,2	0.37	0
5	EDO	B	737	-	3,3,3	0.46	0	2,2,2	0.47	0
5	EDO	A	725	-	3,3,3	0.45	0	2,2,2	0.30	0
5	EDO	B	715	-	3,3,3	0.50	0	2,2,2	0.34	0
5	EDO	A	741	-	3,3,3	0.46	0	2,2,2	0.42	0
5	EDO	B	706	-	3,3,3	0.39	0	2,2,2	0.58	0
5	EDO	A	713	-	3,3,3	0.47	0	2,2,2	0.43	0
5	EDO	B	744	-	3,3,3	0.53	0	2,2,2	0.23	0
5	EDO	B	747	-	3,3,3	0.49	0	2,2,2	0.29	0
5	EDO	A	721	-	3,3,3	0.56	0	2,2,2	0.28	0
5	EDO	A	745	-	3,3,3	0.50	0	2,2,2	0.38	0
6	PGE	A	750	-	9,9,9	0.37	0	8,8,8	0.33	0
5	EDO	A	720	-	3,3,3	0.50	0	2,2,2	0.32	0
5	EDO	B	733	-	3,3,3	0.42	0	2,2,2	0.43	0
5	EDO	B	749[A]	-	3,3,3	0.43	0	2,2,2	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	743	-	3,3,3	0.50	0	2,2,2	0.34	0
5	EDO	A	726	-	3,3,3	0.59	0	2,2,2	0.46	0
5	EDO	B	721	-	3,3,3	0.49	0	2,2,2	0.43	0
5	EDO	A	742	-	3,3,3	0.48	0	2,2,2	0.18	0
5	EDO	A	722	-	3,3,3	0.32	0	2,2,2	0.61	0
5	EDO	A	716	-	3,3,3	0.46	0	2,2,2	0.58	0
5	EDO	B	711	-	3,3,3	0.44	0	2,2,2	0.49	0
5	EDO	B	735	-	3,3,3	0.58	0	2,2,2	0.32	0
5	EDO	B	709	-	3,3,3	0.42	0	2,2,2	0.53	0
5	EDO	B	742	-	3,3,3	0.38	0	2,2,2	0.65	0
5	EDO	B	730	-	3,3,3	0.41	0	2,2,2	0.68	0
5	EDO	A	712	-	3,3,3	0.51	0	2,2,2	0.35	0
5	EDO	B	718	-	3,3,3	0.55	0	2,2,2	0.39	0
5	EDO	B	708	-	3,3,3	0.52	0	2,2,2	0.31	0
8	TRS	B	752	-	7,7,7	0.26	0	9,9,9	0.32	0
5	EDO	B	720	-	3,3,3	0.49	0	2,2,2	0.23	0
5	EDO	A	711	-	3,3,3	0.44	0	2,2,2	0.24	0
7	PEG	B	751	-	6,6,6	0.22	0	5,5,5	0.08	0
5	EDO	B	750	-	3,3,3	0.50	0	2,2,2	0.29	0
5	EDO	A	709	-	3,3,3	0.49	0	2,2,2	0.24	0
5	EDO	A	734	-	3,3,3	0.57	0	2,2,2	0.18	0
5	EDO	B	719	-	3,3,3	0.44	0	2,2,2	0.49	0
5	EDO	A	728	-	3,3,3	0.51	0	2,2,2	0.41	0
5	EDO	A	717	-	3,3,3	0.45	0	2,2,2	0.60	0
5	EDO	A	740	-	3,3,3	0.49	0	2,2,2	0.37	0
5	EDO	B	729	-	3,3,3	0.40	0	2,2,2	0.68	0
5	EDO	A	746	-	3,3,3	0.58	0	2,2,2	0.12	0
5	EDO	A	738	-	3,3,3	0.50	0	2,2,2	0.34	0
5	EDO	A	737	-	3,3,3	0.48	0	2,2,2	0.36	0
5	EDO	B	734	-	3,3,3	0.46	0	2,2,2	0.46	0
5	EDO	A	729	-	3,3,3	0.50	0	2,2,2	0.33	0
5	EDO	B	713	-	3,3,3	0.51	0	2,2,2	0.27	0
5	EDO	A	719	-	3,3,3	0.49	0	2,2,2	0.26	0
5	EDO	A	710	-	3,3,3	0.50	0	2,2,2	0.32	0
5	EDO	B	741	-	3,3,3	0.47	0	2,2,2	0.49	0
5	EDO	B	716	-	3,3,3	0.47	0	2,2,2	0.40	0
5	EDO	B	739	-	3,3,3	0.47	0	2,2,2	0.31	0
5	EDO	A	727	-	3,3,3	0.52	0	2,2,2	0.30	0
5	EDO	B	728	-	3,3,3	0.51	0	2,2,2	0.48	0
5	EDO	B	745	-	3,3,3	0.47	0	2,2,2	0.40	0
5	EDO	A	715	-	3,3,3	0.41	0	2,2,2	0.51	0
5	EDO	A	744	-	3,3,3	0.47	0	2,2,2	0.34	0
5	EDO	A	736	-	3,3,3	0.53	0	2,2,2	0.19	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	EDO	B	722	-	3,3,3	0.49	0	2,2,2	0.28	0
5	EDO	B	736	-	3,3,3	0.50	0	2,2,2	0.26	0
5	EDO	A	743	-	3,3,3	0.41	0	2,2,2	0.53	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	732	-	-	0/1/1/1	-
5	EDO	A	723	-	-	1/1/1/1	-
5	EDO	A	747	-	-	1/1/1/1	-
5	EDO	A	724	-	-	0/1/1/1	-
5	EDO	B	740	-	-	1/1/1/1	-
5	EDO	B	712	-	-	0/1/1/1	-
5	EDO	B	738	-	-	1/1/1/1	-
5	EDO	A	707	-	-	0/1/1/1	-
5	EDO	A	708	-	-	0/1/1/1	-
5	EDO	B	731	-	-	0/1/1/1	-
5	EDO	A	735	-	-	1/1/1/1	-
5	EDO	B	732	-	-	0/1/1/1	-
5	EDO	B	748	-	-	1/1/1/1	-
5	EDO	B	724	-	-	1/1/1/1	-
5	EDO	A	748	-	-	1/1/1/1	-
5	EDO	B	717	-	-	0/1/1/1	-
5	EDO	B	726	-	-	0/1/1/1	-
5	EDO	B	707	-	-	0/1/1/1	-
5	EDO	A	718	-	-	0/1/1/1	-
5	EDO	B	746	-	-	0/1/1/1	-
5	EDO	A	739	-	-	0/1/1/1	-
5	EDO	B	725	-	-	0/1/1/1	-
5	EDO	A	714	-	-	0/1/1/1	-
5	EDO	A	733	-	-	0/1/1/1	-
5	EDO	A	749	-	-	1/1/1/1	-
5	EDO	A	706	-	-	0/1/1/1	-
5	EDO	A	731	-	-	1/1/1/1	-
5	EDO	B	723	-	-	1/1/1/1	-
5	EDO	B	710	-	-	0/1/1/1	-
5	EDO	B	714	-	-	1/1/1/1	-
5	EDO	B	727	-	-	0/1/1/1	-
5	EDO	A	730	-	-	0/1/1/1	-
5	EDO	B	737	-	-	0/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	A	725	-	-	0/1/1/1	-
5	EDO	B	715	-	-	1/1/1/1	-
5	EDO	A	741	-	-	1/1/1/1	-
5	EDO	B	706	-	-	0/1/1/1	-
5	EDO	A	713	-	-	1/1/1/1	-
5	EDO	B	744	-	-	1/1/1/1	-
5	EDO	B	747	-	-	0/1/1/1	-
5	EDO	A	721	-	-	0/1/1/1	-
5	EDO	A	745	-	-	1/1/1/1	-
6	PGE	A	750	-	-	3/7/7/7	-
5	EDO	A	720	-	-	1/1/1/1	-
5	EDO	B	733	-	-	1/1/1/1	-
5	EDO	B	749[A]	-	-	1/1/1/1	-
5	EDO	B	743	-	-	0/1/1/1	-
5	EDO	A	726	-	-	1/1/1/1	-
5	EDO	B	721	-	-	0/1/1/1	-
5	EDO	A	742	-	-	1/1/1/1	-
5	EDO	A	722	-	-	0/1/1/1	-
5	EDO	A	716	-	-	0/1/1/1	-
5	EDO	B	711	-	-	1/1/1/1	-
5	EDO	B	735	-	-	1/1/1/1	-
5	EDO	B	709	-	-	1/1/1/1	-
5	EDO	B	742	-	-	1/1/1/1	-
5	EDO	B	730	-	-	0/1/1/1	-
5	EDO	A	712	-	-	0/1/1/1	-
5	EDO	B	718	-	-	1/1/1/1	-
5	EDO	B	708	-	-	0/1/1/1	-
8	TRS	B	752	-	-	6/9/9/9	-
5	EDO	B	720	-	-	1/1/1/1	-
5	EDO	A	711	-	-	0/1/1/1	-
7	PEG	B	751	-	-	3/4/4/4	-
5	EDO	B	750	-	-	0/1/1/1	-
5	EDO	A	709	-	-	1/1/1/1	-
5	EDO	A	734	-	-	1/1/1/1	-
5	EDO	B	719	-	-	0/1/1/1	-
5	EDO	A	728	-	-	1/1/1/1	-
5	EDO	A	717	-	-	0/1/1/1	-
5	EDO	A	740	-	-	0/1/1/1	-
5	EDO	B	729	-	-	1/1/1/1	-
5	EDO	A	746	-	-	1/1/1/1	-
5	EDO	A	738	-	-	1/1/1/1	-
5	EDO	A	737	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	B	734	-	-	1/1/1/1	-
5	EDO	A	729	-	-	0/1/1/1	-
5	EDO	B	713	-	-	0/1/1/1	-
5	EDO	A	719	-	-	1/1/1/1	-
5	EDO	A	710	-	-	1/1/1/1	-
5	EDO	B	741	-	-	0/1/1/1	-
5	EDO	B	716	-	-	0/1/1/1	-
5	EDO	B	739	-	-	0/1/1/1	-
5	EDO	A	727	-	-	0/1/1/1	-
5	EDO	B	728	-	-	0/1/1/1	-
5	EDO	B	745	-	-	1/1/1/1	-
5	EDO	A	715	-	-	1/1/1/1	-
5	EDO	A	744	-	-	0/1/1/1	-
5	EDO	A	736	-	-	0/1/1/1	-
5	EDO	B	722	-	-	0/1/1/1	-
5	EDO	B	736	-	-	0/1/1/1	-
5	EDO	A	743	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 53 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	752	TRS	C1-C-C2-O2
8	B	752	TRS	C3-C-C2-O2
7	B	751	PEG	O2-C3-C4-O4
5	A	709	EDO	O1-C1-C2-O2
5	A	710	EDO	O1-C1-C2-O2

There are no ring outliers.

33 monomers are involved in 49 short contacts:

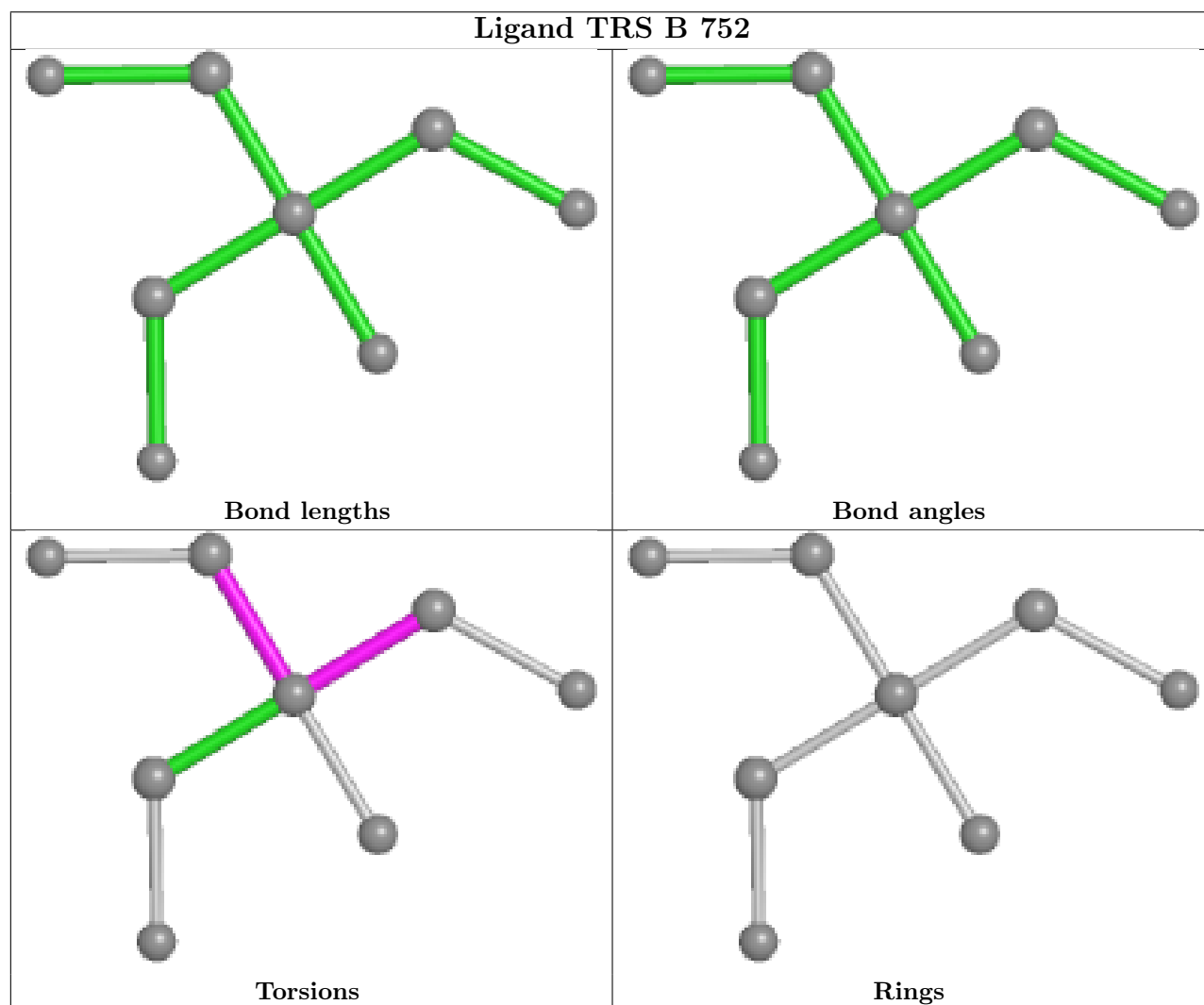
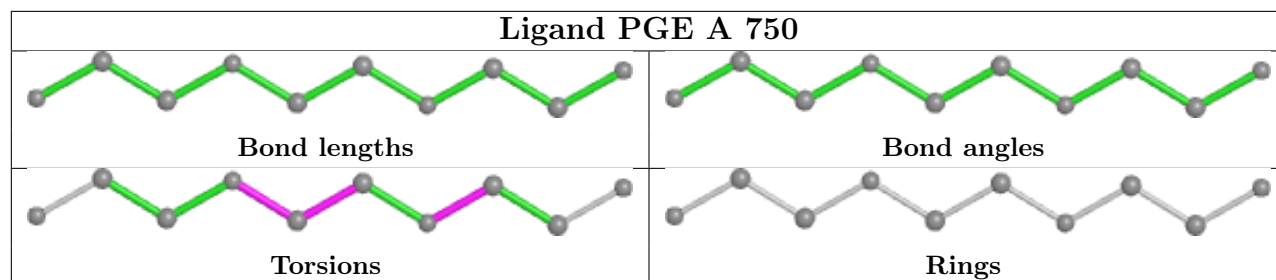
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	723	EDO	1	0
5	A	724	EDO	3	0
5	B	732	EDO	4	0
5	B	724	EDO	1	0
5	B	746	EDO	1	0
5	A	739	EDO	2	0
5	B	723	EDO	1	0

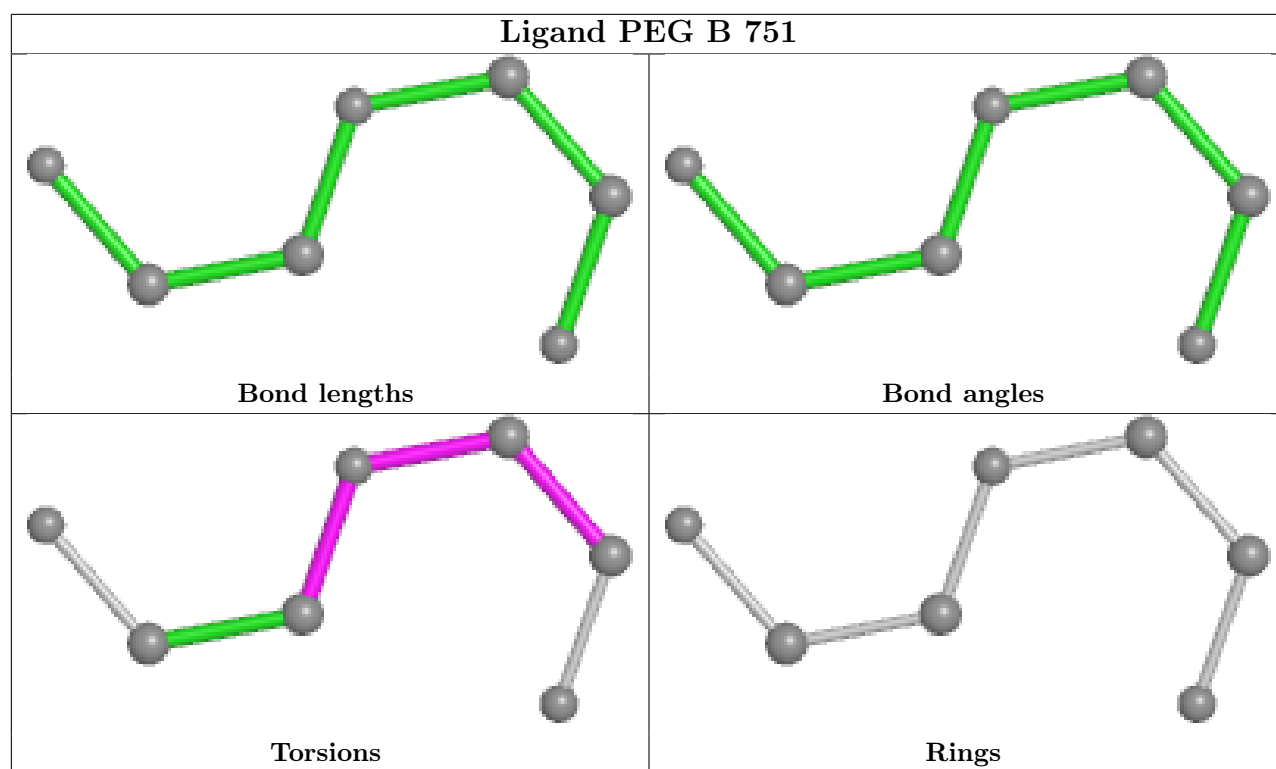
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	714	EDO	1	0
5	B	727	EDO	1	0
5	A	730	EDO	1	0
5	B	737	EDO	1	0
5	A	725	EDO	2	0
5	A	741	EDO	1	0
5	A	713	EDO	1	0
5	B	743	EDO	1	0
5	A	726	EDO	1	0
5	A	722	EDO	3	0
5	B	735	EDO	3	0
5	B	709	EDO	1	0
5	B	742	EDO	4	0
5	B	730	EDO	1	0
5	B	718	EDO	1	0
7	B	751	PEG	2	0
5	A	728	EDO	2	0
5	B	729	EDO	1	0
5	A	737	EDO	1	0
5	B	734	EDO	1	0
5	A	719	EDO	1	0
5	A	710	EDO	1	0
5	B	716	EDO	1	0
5	A	727	EDO	1	0
5	A	744	EDO	1	0
5	B	736	EDO	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	532/559 (95%)	-0.17	8 (1%) 73 72	21, 31, 55, 86	0
1	B	532/559 (95%)	-0.35	3 (0%) 89 88	20, 29, 48, 77	0
2	C	4/6 (66%)	1.19	1 (25%) 0 0	34, 36, 39, 40	4 (100%)
2	E	4/6 (66%)	0.06	0 100 100	36, 38, 42, 48	0
2	F	2/6 (33%)	0.63	0 100 100	45, 45, 45, 50	2 (100%)
2	G	4/6 (66%)	1.09	1 (25%) 0 0	34, 34, 37, 40	4 (100%)
2	I	4/6 (66%)	-0.57	0 100 100	34, 36, 39, 40	0
2	J	2/6 (33%)	2.67	1 (50%) 0 0	53, 53, 53, 56	0
All	All	1084/1154 (93%)	-0.24	14 (1%) 77 76	20, 30, 51, 86	10 (0%)

The worst 5 of 14 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	J	4	GLY	3.8
1	A	92	CYS	3.7
1	A	130	ALA	3.1
1	A	94	LEU	3.0
1	B	362	ASN	2.8

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	HYP	G	6	9/9	0.80	0.15	34,37,41,42	9

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	HYP	C	6	9/9	0.85	0.15	35,38,41,42	9
2	HYP	J	6	8/9	0.85	0.13	35,40,49,49	0
2	HYP	F	6	8/9	0.87	0.14	33,37,47,47	8
2	HYP	C	3	8/9	0.88	0.15	35,38,43,45	8
2	HYP	G	3	8/9	0.90	0.12	34,37,40,42	8
2	HYP	E	6	9/9	0.91	0.12	36,42,48,61	0
2	HYP	I	6	9/9	0.93	0.11	34,35,45,55	0
2	HYP	I	3	8/9	0.96	0.10	27,34,37,44	0
2	HYP	E	3	8/9	0.98	0.10	31,34,35,40	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	EDO	B	714	4/4	0.60	0.24	46,50,50,55	0
5	EDO	A	721	4/4	0.61	0.20	45,49,50,54	0
5	EDO	B	743	4/4	0.61	0.21	53,55,57,62	0
5	EDO	B	740	4/4	0.70	0.19	46,47,48,57	0
5	EDO	A	746	4/4	0.71	0.22	52,55,56,61	0
5	EDO	A	735	4/4	0.72	0.20	43,46,47,47	0
8	TRS	B	752	8/8	0.74	0.18	54,56,64,65	0
5	EDO	A	730	4/4	0.76	0.25	48,50,51,57	0
7	PEG	B	751	7/7	0.77	0.21	30,47,52,54	0
5	EDO	A	728	4/4	0.77	0.30	42,46,46,48	0
5	EDO	B	722	4/4	0.78	0.23	65,66,69,72	0
5	EDO	A	733	4/4	0.79	0.18	48,52,56,56	0
6	PGE	A	750	10/10	0.79	0.23	38,48,56,59	0
5	EDO	A	718	4/4	0.80	0.14	46,47,48,51	0
5	EDO	A	736	4/4	0.80	0.12	54,58,60,61	0
5	EDO	B	726	4/4	0.80	0.15	51,51,54,64	0
5	EDO	A	720	4/4	0.80	0.27	46,48,50,53	0
5	EDO	B	747	4/4	0.81	0.23	56,57,58,65	0
5	EDO	B	736	4/4	0.81	0.32	49,50,52,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	EDO	B	750	4/4	0.82	0.13	42,44,50,56	0
5	EDO	B	712	4/4	0.82	0.20	47,52,53,53	0
5	EDO	A	734	4/4	0.82	0.20	39,40,44,52	0
5	EDO	B	717	4/4	0.82	0.15	41,41,43,46	0
5	EDO	A	749	4/4	0.83	0.16	34,48,53,62	0
5	EDO	B	749[A]	4/4	0.83	0.16	26,33,36,39	4
5	EDO	B	716	4/4	0.83	0.14	49,49,49,55	0
5	EDO	A	731	4/4	0.83	0.16	33,39,43,52	0
5	EDO	B	718	4/4	0.83	0.20	36,48,50,52	0
5	EDO	B	721	4/4	0.83	0.30	52,55,55,61	0
5	EDO	A	748	4/4	0.84	0.19	40,46,47,58	0
5	EDO	B	725	4/4	0.84	0.16	37,48,52,62	0
5	EDO	B	709	4/4	0.85	0.26	37,51,55,67	0
5	EDO	B	723	4/4	0.86	0.24	46,47,50,54	0
5	EDO	A	726	4/4	0.86	0.17	31,41,43,44	0
5	EDO	A	727	4/4	0.86	0.14	27,39,49,49	0
5	EDO	A	714	4/4	0.86	0.11	43,44,45,59	0
5	EDO	A	719	4/4	0.86	0.20	35,42,52,54	0
5	EDO	A	725	4/4	0.86	0.19	39,46,47,52	0
5	EDO	A	745	4/4	0.87	0.27	45,52,53,54	0
5	EDO	B	710	4/4	0.87	0.22	31,37,44,46	0
5	EDO	A	707	4/4	0.87	0.17	47,48,51,54	0
5	EDO	B	729	4/4	0.87	0.22	38,39,44,47	0
5	EDO	B	713	4/4	0.87	0.10	36,40,41,46	0
5	EDO	B	739	4/4	0.87	0.23	48,49,50,57	0
5	EDO	B	707	4/4	0.87	0.19	33,39,45,52	0
5	EDO	A	747	4/4	0.88	0.10	36,44,52,56	0
5	EDO	A	710	4/4	0.88	0.19	37,43,44,46	0
5	EDO	A	712	4/4	0.89	0.18	35,50,53,58	0
5	EDO	A	723	4/4	0.89	0.17	34,34,36,49	0
5	EDO	A	713	4/4	0.89	0.17	29,41,45,46	0
5	EDO	A	740	4/4	0.90	0.10	38,39,46,47	0
5	EDO	A	743	4/4	0.90	0.17	37,38,44,51	0
5	EDO	B	730	4/4	0.90	0.18	44,45,45,46	0
5	EDO	B	735	4/4	0.90	0.28	24,31,41,44	0
5	EDO	A	715	4/4	0.90	0.18	44,45,46,50	0
5	EDO	A	708	4/4	0.90	0.14	40,53,54,57	0
5	EDO	A	738	4/4	0.90	0.23	39,42,46,56	0
5	EDO	B	719	4/4	0.91	0.15	39,48,52,66	0
5	EDO	B	715	4/4	0.91	0.25	38,42,50,51	0
4	CA	A	705	1/1	0.91	0.43	29,29,29,29	0
5	EDO	B	744	4/4	0.91	0.10	39,60,60,64	0

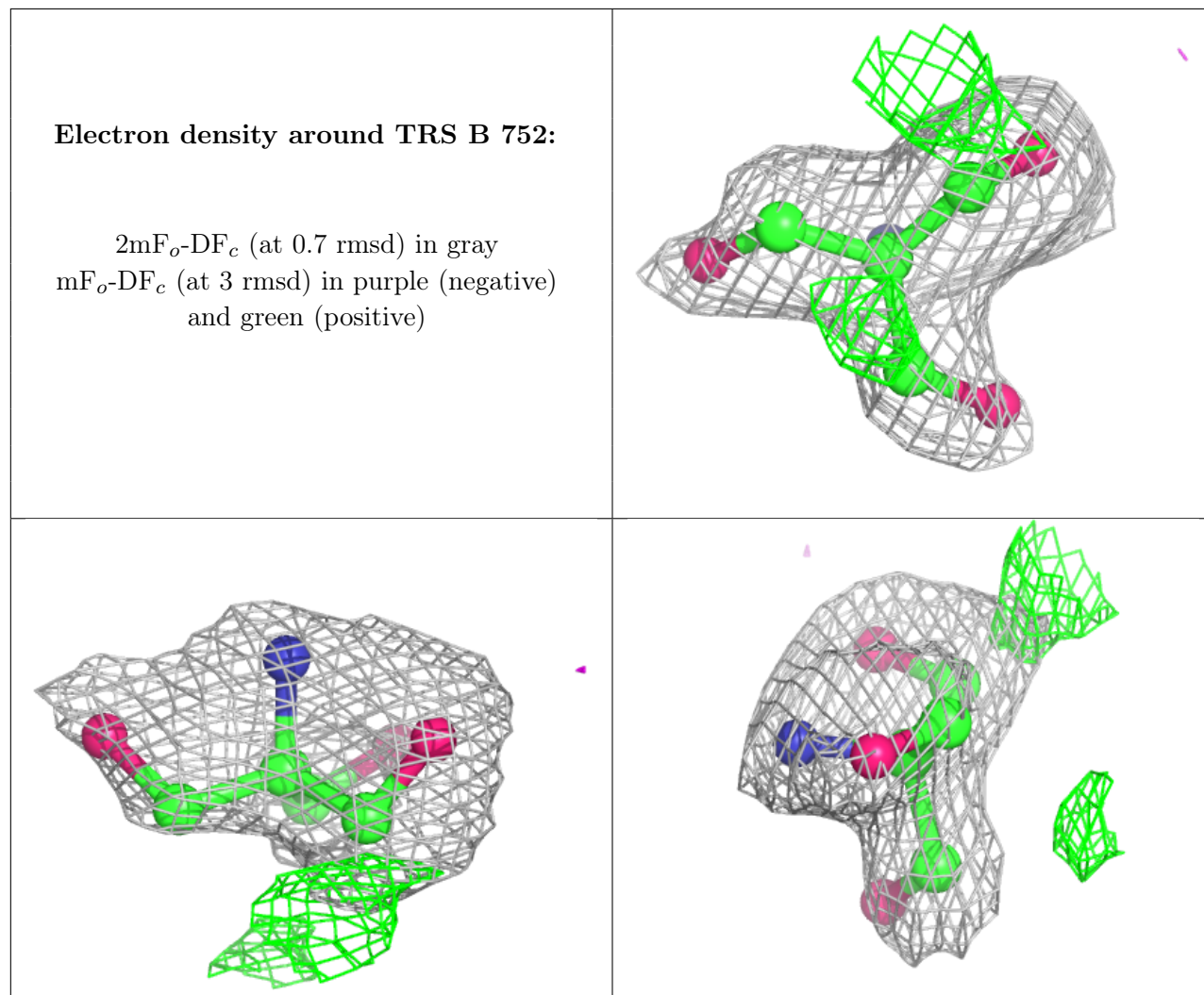
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	EDO	B	727	4/4	0.91	0.12	42,44,49,53	0
5	EDO	A	742	4/4	0.92	0.19	29,37,40,52	0
5	EDO	B	737	4/4	0.92	0.09	38,42,44,45	0
4	CA	B	705	1/1	0.92	0.17	62,62,62,62	0
5	EDO	B	731	4/4	0.92	0.10	43,48,50,60	0
5	EDO	B	733	4/4	0.92	0.16	35,36,39,44	0
5	EDO	A	709	4/4	0.92	0.18	32,37,39,50	0
5	EDO	A	741	4/4	0.93	0.17	36,42,48,52	0
5	EDO	A	729	4/4	0.93	0.15	30,35,36,38	0
5	EDO	B	732	4/4	0.93	0.15	24,31,36,39	0
5	EDO	B	742	4/4	0.93	0.14	31,38,48,60	0
5	EDO	A	737	4/4	0.94	0.11	39,43,44,47	0
5	EDO	A	706	4/4	0.94	0.14	29,33,37,51	0
5	EDO	A	744	4/4	0.94	0.11	54,54,56,64	0
5	EDO	B	746	4/4	0.94	0.17	32,52,55,57	0
5	EDO	B	734	4/4	0.94	0.17	32,36,38,50	0
5	EDO	B	748	4/4	0.94	0.20	38,41,47,59	0
5	EDO	A	739	4/4	0.94	0.33	36,40,41,57	0
5	EDO	A	716	4/4	0.94	0.12	37,39,41,42	0
5	EDO	B	711	4/4	0.94	0.16	32,35,36,38	0
5	EDO	A	722	4/4	0.94	0.27	34,36,37,43	0
5	EDO	B	720	4/4	0.94	0.31	45,54,58,61	0
5	EDO	B	708	4/4	0.95	0.09	30,39,47,47	0
5	EDO	B	728	4/4	0.95	0.11	29,32,34,40	0
5	EDO	B	738	4/4	0.95	0.24	40,42,45,46	0
5	EDO	A	732	4/4	0.95	0.12	36,41,42,45	0
5	EDO	A	711	4/4	0.96	0.12	37,41,44,51	0
5	EDO	B	745	4/4	0.96	0.16	35,41,44,63	0
5	EDO	B	724	4/4	0.96	0.10	33,38,41,45	0
5	EDO	B	706	4/4	0.96	0.13	24,35,39,47	0
5	EDO	A	717	4/4	0.96	0.08	30,38,38,52	0
5	EDO	A	724	4/4	0.97	0.16	30,38,45,49	0
3	ZN	B	701	1/1	0.98	0.06	38,38,38,38	1
5	EDO	B	741	4/4	0.98	0.11	38,39,40,51	0
3	ZN	A	701	1/1	0.99	0.09	40,40,40,40	1
4	CA	B	703	1/1	0.99	0.06	30,30,30,30	0
4	CA	B	704	1/1	0.99	0.03	33,33,33,33	0
4	CA	A	702	1/1	0.99	0.12	21,21,21,21	0
4	CA	A	703	1/1	0.99	0.07	32,32,32,32	0
4	CA	A	704	1/1	0.99	0.04	35,35,35,35	0
4	CA	B	702	1/1	1.00	0.12	23,23,23,23	0

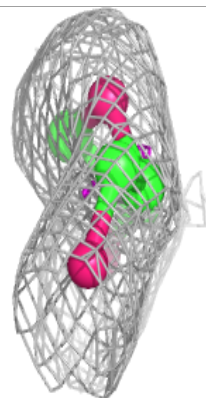
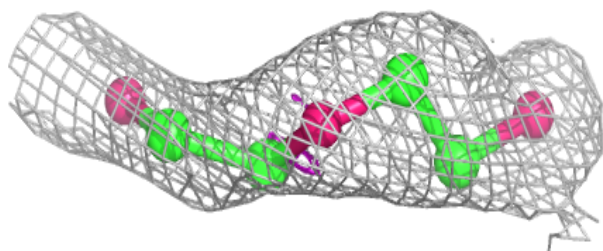
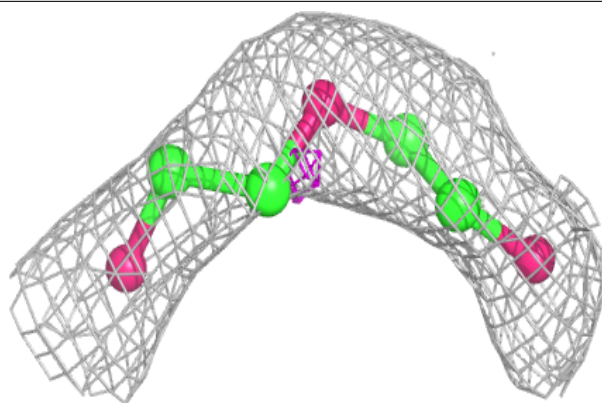
The following is a graphical depiction of the model fit to experimental electron density of all

instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

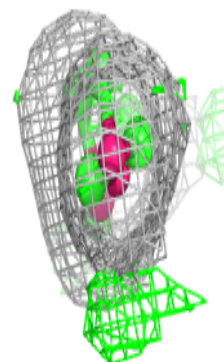
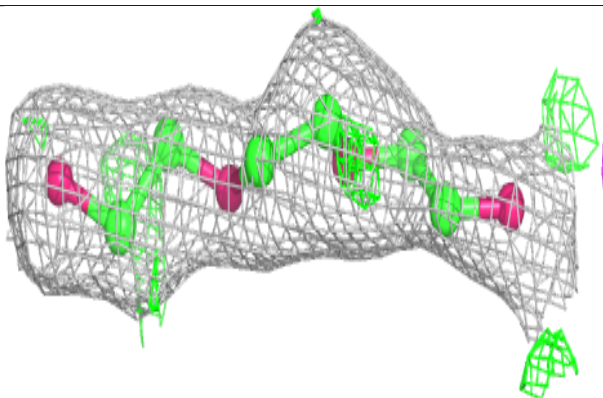
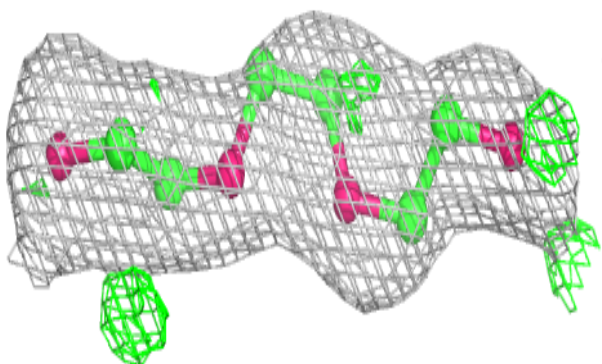


Electron density around PEG B 751:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

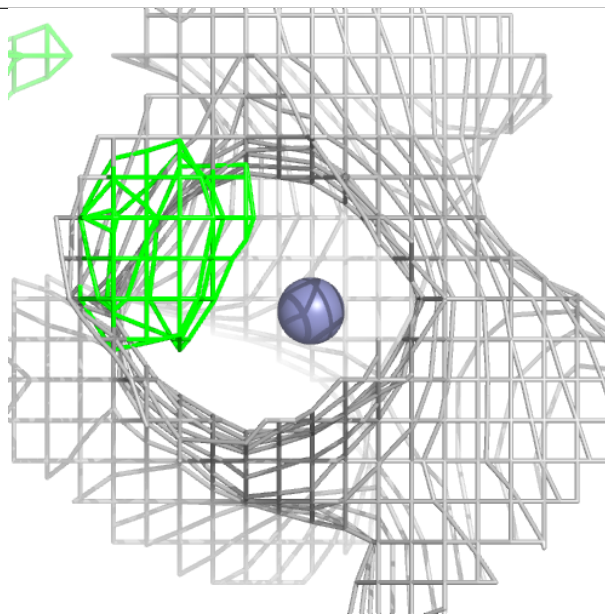
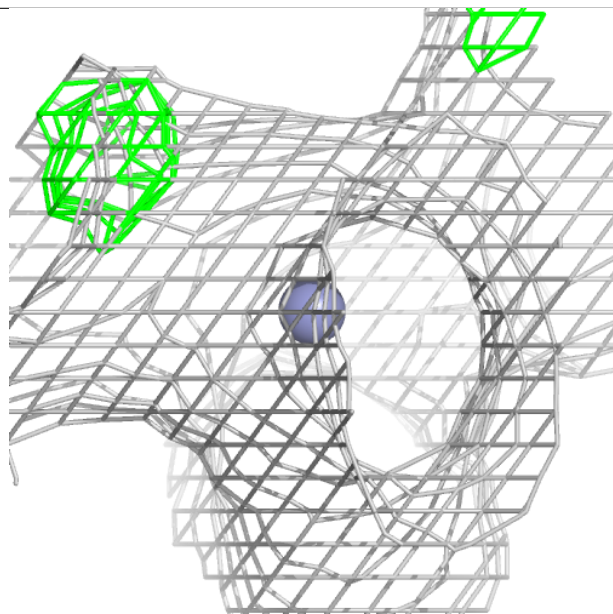
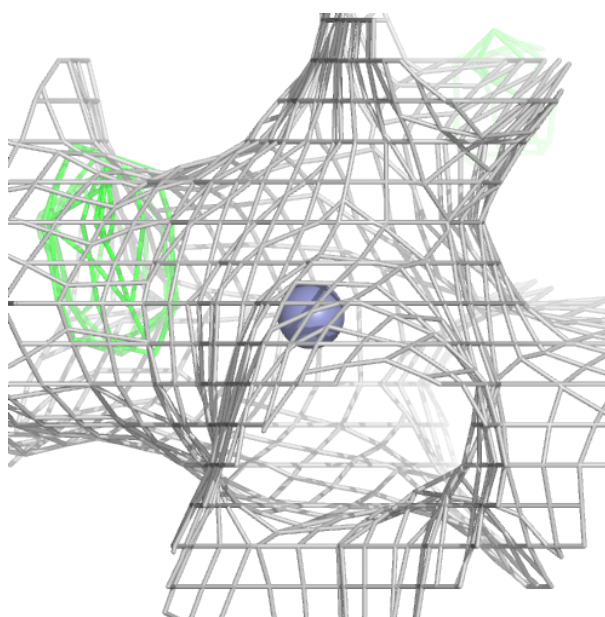
**Electron density around PGE A 750:**

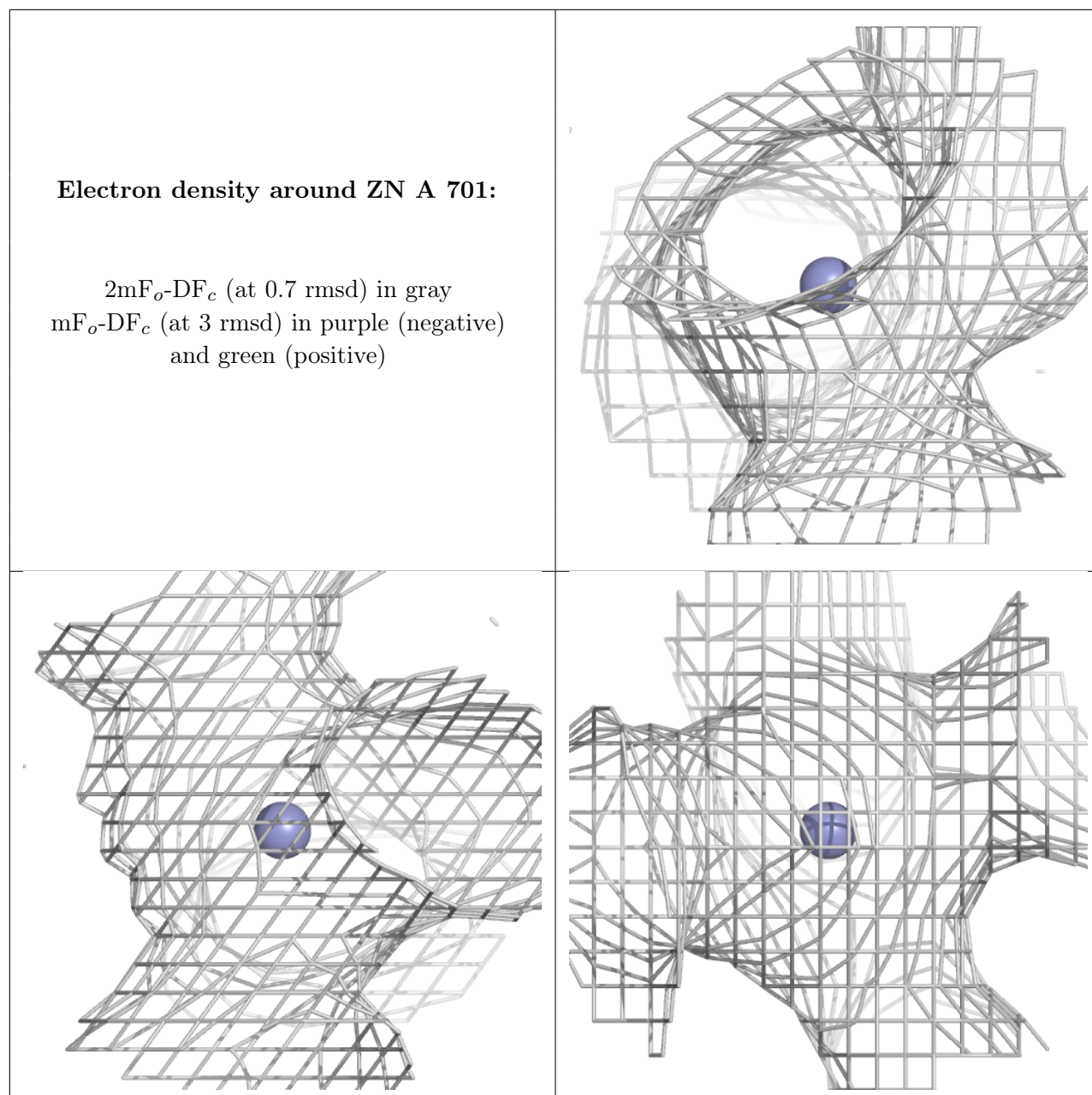
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around ZN B 701:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.